

Data Path : Z:\HPCHEM1\BNA F\DATA\BF101617\
 Data File : BF099569.D
 Acq On : 17 Oct 2017 2:52
 Operator : SJ/JU
 Sample : I5752-10RE
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEST-PIT-116-C(0-5)RE

Manual Integrations
 APPROVED

Sohil
 10/17/2017 5:02:08 PM

Quant Time: Oct 17 03:54:01 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 11 16:25:21 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	99429	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	400421	20.00	ng	0.00
38) Acenaphthene-d10	9.88	164	167863	20.00	ng	-0.01
63) Phenanthrene-d10	11.37	188	247550	20.00	ng	0.00
75) Chrysene-d12	14.02	240	177968	20.00	ng	0.00
86) Perylene-d12	15.49	264	189713	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	704798	112.84	ng	0.00
7) Phenol-d6	6.48	99	838890	108.88	ng	0.00
23) Nitrobenzene-d5	7.41	82	493202	78.99	ng	-0.01
41) 2,4,6-Tribromophenol	10.67	330	144401	105.13	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	854343	84.97	ng	-0.01
78) Terphenyl-d14	12.95	244	535351	68.41	ng	0.00
Target Compounds						
10) Phenol	6.49	94	50887	5.905	ng	97
49) Dimethylphthalate	9.59	163	127774	10.085	ng	98
70) Phenanthrene	11.39	178	54402	4.106	ng	96
74) Fluoranthene	12.59	202	74800	5.575	ng	97
77) Pyrene	12.82	202	68178	5.024	ng	98
80) Benzo(a)anthracene	14.00	228	36216	3.361	ng	97
82) Chrysene	14.04	228	29541	2.879	ng	97
85) Indeno(1,2,3-cd)pyrene	16.97	276	18374	2.239	ng	96
87) Benzo(b)fluoranthene	15.06	252	42170m	3.615	ng	
89) Benzo(a)pyrene	15.43	252	31469	2.939	ng	# 92

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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